

# **Structure and intramolecular lability of N-(thio)phosphoryl(thio)amides: VI. $^1\text{H}$ , $^{13}\text{C}$ , and $^{31}\text{P}$ NMR interpretation of the phosphorylotropic rearrangement in the N,N'-bis(diisopropoxythiophosphorylaminocarbonyl)-1,10-diaza-18-crown-6 ether in DMSO solution**

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## **Abstract**

The phosphorylotropic rearrangement in the N,N'-bis(diisopropoxythiophosphorylaminocarbonyl)-1,10-diaza-18-crown-6 ether in DMSO solution was studied by  $^1\text{H}$ ,  $^{13}\text{C}$ , and  $^{31}\text{P}$  NMR spectroscopy. The transition from the amide to phosphorylotropic form is accompanied by simultaneous transformation of signals of all structural elements in the  $^1\text{H}$ ,  $^{13}\text{C}$ , and  $^{31}\text{P}$  NMR spectra.

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